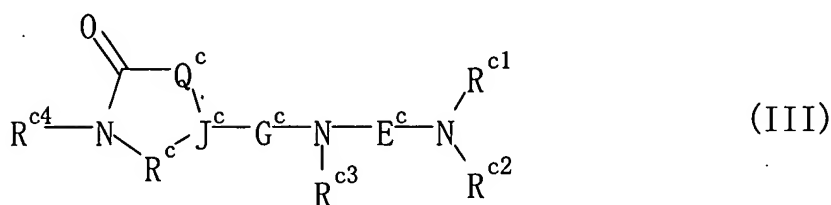


$$\begin{array}{c}
 \text{R}^{a4}-\text{G}^{a1}-\text{N} \begin{array}{l} \nearrow \text{Q}^a \\ \searrow \text{R}^a \end{array} \text{J}^a-\text{G}^{a2}-\text{N} \begin{array}{l} \nearrow \text{E}^a \\ \searrow \text{A}^a \end{array} \begin{array}{l} \text{R}^{a1} \\ \text{R}^{a2} \end{array} \\
 \text{R}^{a3} \\
 \text{(CH}_2\text{)}_{na}
 \end{array} \quad (I)$$
$$\begin{array}{c} \text{O} \\ || \\ R^{b1}-\text{NH}-\text{N}-\text{CH}_2-\text{CH}_2-\text{N}-\text{C}_6\text{H}_4-(R^{b3})_{pb} \\ | \\ (\text{CH}_2)_{nb} \\ | \\ R^{b2} \end{array}$$

(II)

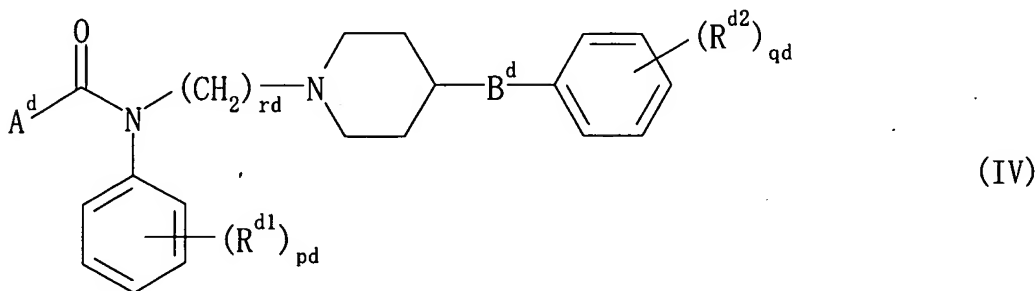
wherein R^{b1} is a hydrocarbon group which may be substituted; R^{b2} is a cyclic hydrocarbon group which may be substituted or a heterocyclic group which may be substituted; R^{b3} is a halogen atom, a carbamoyl group which may be substituted, a sulfamoyl group which may be substituted, an acyl group derived from a sulfonic acid, a C_{1-4} alkyl group which may be substituted, a C_{1-4} alkoxy group which may be substituted, an amino group which may be substituted, a nitro group or a cyano group; R^{b4} is a hydrogen atom or a hydroxy group; nb is an integer of 0 or 1; pb is an integer of 0 or 1 to 4,

the formula:

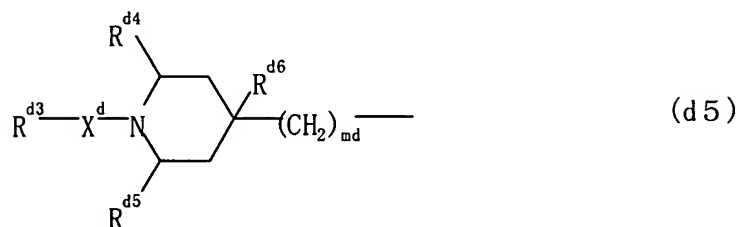
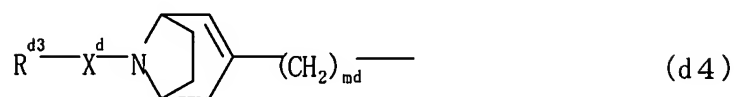
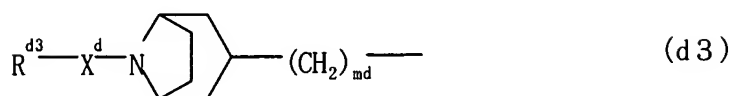
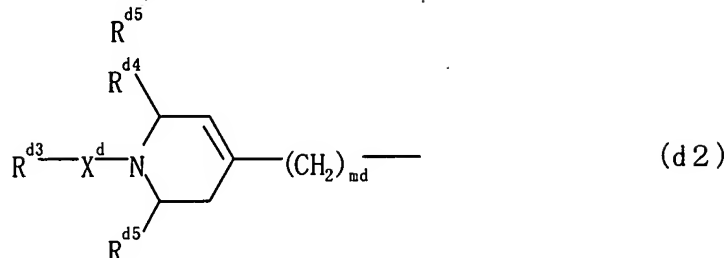
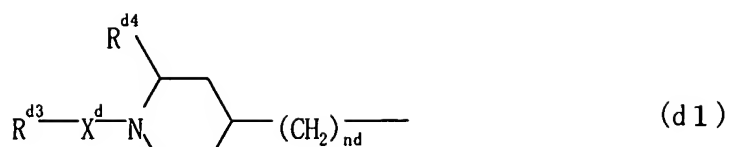


wherein R^{c1} is a hydrocarbon group, R^{c2} is a hydrocarbon group having 2 or more carbon atoms, or R^{c1} and R^{c2} may be bound together with the adjacent nitrogen atom to form a ring which may have a substituent or substituents, R^{c3} is a hydrocarbon group which may have a substituent or substituents or a heterocyclic group which may have a substituent or substituents, R^{c4} is a hydrogen atom, a hydrocarbon group which may have a substituent or substituents or a heterocyclic group which may have a substituent or substituents, E^c is a divalent aliphatic hydrocarbon group which may have a substituent or substituents other than an oxo group, G^c is CO or SO₂, J^c is a nitrogen atom or a methine group which may have a substituent or substituents, and Q^c and R^c are each a bond or a divalent aliphatic C_{1-3} hydrocarbon group which may have a substituent or substituents,

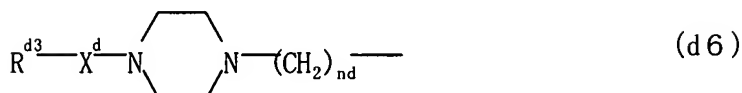
the formula:



wherein A^d is a group represented by the formula:

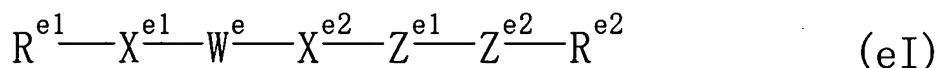


or

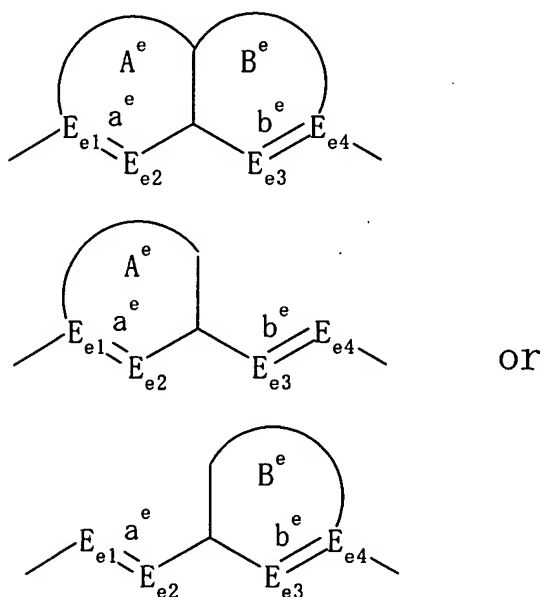


wherein, R^{d3} is (1) a hydrocarbon group which may be substituted, (2) a C_{1-4} alkoxy group which may be substituted or (3) an amino group which may be substituted; X^d is a bond, $-SO_2-$ or $-CO-$; nd is an integer of 1 to 3; md is 0 or an integer of 1 to 3; R^{d4} and R^{d5} are the same or different and each of which is a hydrogen atom or a C_{1-6} alkyl group; R^{d6} is a hydroxyl group, a C_{1-6} alkyl group or a C_{2-6} alkenyl group; rd is an integer of 2 to 4; B^d is a bond, $-CH_2-$, $-SO_2-$, $-SO-$, $-S-$, $-O-$, $-CO-$, $-NR^{da}-SO_2-$ or $-NR^{da}-CO-$ (wherein, R^{da} is a hydrogen atom, a C_{1-6} alkyl group, a C_{2-6} alkenyl group or a C_{3-8} cycloalkyl group); each of pd and qd is 0 or an integer of 1 to 4; R^{d1} is a halogen atom, a C_{1-6} alkyl group, a C_{2-4} alkenyl group, a C_{1-4} alkanoyl group, a C_{1-4} alkoxy group, a cyano group, a trifluoromethyl group, a nitro group, a hydroxyl group, an amino group or an amidino group; R^{d2} is 1) a halogen, 2) a C_{1-6} alkyl which may be substituted by a halogen or a

C₁₋₄ alkoxy, 3) a C₁₋₄ alkoxy which may be substituted by a halogen or a C₁₋₄ alkoxy, 4) nitro, 5) cyano, 6) hydroxyl, 7) a C₁₋₄ alkanoylamino, 8) SO₂NR^{db}R^{dc}, 9) SO₂R^{dd}, 10) CONR^{db}R^{dc}, 11) NR^{db}R^{dc} or 12) NR^{da}-SO₂R^{dd} (wherein, R^{da} has the meaning given above, and R^{db} and R^{dc} may be the same or different, and are (1) a hydrogen atom, (2) a C₁₋₆ alkyl group which may be substituted by a halogen or a C₁₋₄ alkoxy, or (3) a C₃₋₈ cycloalkyl group which may be substituted by a halogen or a C₁₋₄ alkoxy, or R^{db} and R^{dc} may bond with a nitrogen atom to form a cyclic amino group and R^{dd} is a C₁₋₆ alkyl group or a C₃₋₈ cycloalkyl group), each R^{d1} may be the same or different from each other when pd is two or more, and each R^{d2}, may be the same or different from each other when qd is two or more, or the formula:

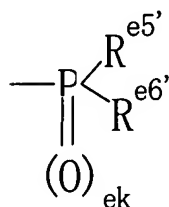


wherein R^{e1} represents a 5 to 6-membered cyclic ring group which may be substituted, X^{e1} represents a bond or a bivalent group, in which the number of atoms constituting the straight-chain portion is 1 to 4, W^e represents a bivalent group represented by formula:



wherein each of ring A^e and ring B^e represents a 5- to 7-membered cyclic group which may be substituted, each of E_{e1} and E_{e4} is a carbon atom which may be substituted or a nitrogen atom which may be substituted, each of E_{e2} and E_{e3} is a carbon atom which may be substituted, a nitrogen atom which may be substituted, or a sulfur atom which may be oxidized or an oxygen

atom, each of a^e and b^e is a single bond or a double bond), X^{e2} is a bivalent group in which the number of atoms constituting the straight-chain portion is 1 to 4, Z^{e1} is a bond or a bivalent cyclic ring group, Z^{e2} is a bond or a bivalent cyclic ring group in which the number of atoms constituting the straight-chain portion is 1 to 4, and R^{e2} is (1) an amino group which may be substituted, and the nitrogen atom may be converted into a quaternary ammonium or an N-oxide, (2) a nitrogen-containing heterocyclic ring group which may be substituted, may contain sulfur atom or an oxygen atom as a ring-constituting atom, and the nitrogen atom may be converted into a quaternary ammonium or a N-oxide, (3) a group which is bonded via the sulfur atom, (4) a group represented by formula:



wherein ek is 0 or 1, the phosphorus atom may form a phosphonium salt when ek is 0, and each of $R^{e5'}$ and $R^{e6'}$ is a hydrocarbon atom which may be substituted, a hydroxyl group which may be substituted, or an amino group which may be substituted, and $R^{e5'}$ and $R^{e6'}$ may bond to each other to form a cyclic ring group together with the adjacent phosphorus atom, (5) an amidino group which may be substituted or (6) a guanidino group which may be substituted, or a salt thereof.

2. (Cancelled)

3. (Currently amended) The method according to ~~claim 5~~claim 1, wherein the compound having a CCR antagonistic effect or a salt thereof is N-(3,4-dichlorophenyl)-1-(methylsulfonyl)-N-{3-[4-(4-[(methylsulfonyl)-amino]phenyl)sulfonyl]-1-piperidinyl}propyl}-4-piperidinecarboxamide, N-(3-chlorophenyl)-1-(methylsulfonyl)-N-(3-{4-[4-(methylsulfonyl)benzyl]-1-piperidinyl}propyl)-4-piperidinecarboxamide, N-(3-{4-[4-(aminocarbonyl)benzyl]-1-piperidinyl}propyl)-N-(3,4-dichlorophenyl)-1-(methylsulfonyl)-4-piperidinecarboxamide, 1-acetyl-N-(3-{4-[4-(aminocarbonyl)benzyl]-1-piperidinyl}propyl)-N-(3-chloro-4-methylphenyl)-4-piperidinecarboxamide, N-(3,4-dichlorophenyl)-N-(3-{4-[4-

(ethylsulfonyl)benzyl]-1-piperidinyl}propyl)-1-(methylsulfonyl)-4-piperidinecarboxamide, N-(3,4-dichlorophenyl)-N-(3-{4-[4-(isopropylsulfonyl)benzyl]-1-piperidinyl}propyl)-1-(methylsulfonyl)-4-piperidinecarboxamide, N-(3-chlorophenyl)-N-(3-{4-[4-(isopropylsulfonyl)benzyl]-1-piperidinyl}propyl)-1-(methylsulfonyl)-4-piperidinecarboxamide, N-(3-chlorophenyl)-N-(3-{4-[4-(ethylsulfonyl)benzyl]-1-piperidinyl}propyl)-1-(methylsulfonyl)-4-piperidinecarboxamide, N-(3,4-dichlorophenyl)-1-(methylsulfonyl)-N-(3-{4-[4-(methylsulfonyl)benzyl]-1-piperidinyl}propyl)-4-piperidinecarboxamide, N-(3-{4-[4-(aminocarbonyl)benzyl]-1-piperidinyl}propyl)-N-(3-chloro-4-methylphenyl)-1-(methylsulfonyl)-4-piperidinecarboxamide, N-[3-(4-benzyl-1-piperidinyl)propyl]-N'-(4-chlorophenyl)-N-phenylurea, N'-(4-chlorophenyl)-N-{3-[4-(4-fluorobenzyl)-1-piperidinyl]propyl}-N-phenylurea, N'-(4-chlorophenyl)-N-(3-{4-[4-(4-morpholinylsulfonyl)benzyl]-1-piperidinyl}propyl)-N-phenylurea, N'-(4-chlorophenyl)-N-(3-{4-[4-(4-methylsulfonyl)benzyl]-1-piperidinyl}propyl)-N-phenylurea, 4-{[1-(3-{[(4-chloroanilino)carbonyl]anilino}propyl)-4-piperidinyl]methyl}benzamide, N-[3-(4-benzyl-1-piperidinyl)propyl]-N-(3,4-dichlorophenyl)-1-methyl-5-oxo-3-pyrrolidinecarboxamide, 1-benzyl-N-[3-(4-benzyl-1-piperidinyl)propyl]-5-oxo-N-phenyl-3-pyrrolidinecarboxamide, N-[3-(4-benzyl-1-piperidinyl)propyl]-1-(2-chlorobenzyl)-5-oxo-N-phenyl-3-pyrrolidinecarboxamide, N-(3,4-dichlorophenyl)-N-{3-[4-(4-fluorobenzyl)-1-piperidinyl]propyl}-1-methyl-5-oxo-3-pyrrolidinecarboxamide, N-[3-(4-benzyl-1-piperidinyl)propyl]-5-oxo-N-phenyl-1-(2,2,2-trifluoroethyl)-3-pyrrolidinecarboxamide, N-(3,4-dichlorophenyl)-N-(3-{4-[4-(methylsulfonyl)benzyl]-1-piperidinyl}propyl)-2-[1-(methylsulfonyl)-4-piperidinyl]acetamide, N-(3,4-dichlorophenyl)-N-(3-{4-[4-(isopropylsulfonyl)benzyl]-1-piperidinyl}propyl)-2-[1-(methylsulfonyl)-4-piperidinyl]acetamide, 3-(1-acetyl-4-piperidinyl)-N-(3,4-dichlorophenyl)-N-(3-{4-[4-(methylsulfonyl)benzyl]-1-piperidinyl}propyl)propanamide, or N-(3,4-dichlorophenyl)-4-hydroxy-1-(methylsulfonyl)-N-(3-{4-[4-(methylsulfonyl)benzyl]-1-piperidinyl}propyl)-4-piperidinecarboxamide or a salt thereof.

4. (Currently amended) The method according to ~~claim 5~~claim 1, wherein the compound having a CCR antagonistic effect or a salt thereof is N-methyl-N-[4-[[[2-(4-methylphenyl)-6,7-dihydro-5H-benzocyclohepten-8-yl]carbonyl]amino]benzyl]piperidinium iodide, N-methyl-N-[4-[[[7-(4-methylphenyl)-2,3-dihydro-1-benzoxepin-4-yl]carbonyl]amino]benzyl]piperidinium

iodide, N-[4-[N-methyl-N-(tetrahydropyran-4-yl)aminomethyl]phenyl]-7-(4-methylphenyl)-2,3-dihydro-1-benzoxepin-4-carboxamide, N-[4-[N-methyl-N-(tetrahydropyran-4-yl)aminomethyl]phenyl]-7-(4-morpholinophenyl)-2,3-dihydro-1-benzoxepin-4-carboxamide, 7-(4-ethoxyphenyl)-N-[4-[N-methyl-N-(tetrahydropyran-4-yl)aminomethyl]phenyl]-2,3-dihydro-1-benzoxepin-4-carboxamide, N,N-dimethyl-N-[4-[[[2-(4-methylphenyl)-6,7-dihydro-5H-benzocyclohepten-8-yl]carbonyl]amino]benzyl]-N-(tetrahydropyran-4-yl)ammonium iodide, N-methyl-N-[4-[[[7-(4-methylphenyl)-3,4-dihydronaphthalen-2-yl]carbonyl]amino]benzyl]piperidinium iodide, N,N-dimethyl-N-(4-(((2-(4-methylphenyl)-6,7-dihydro-5H-benzocyclohepten-8-yl)carbonyl)amino)benzyl)-N-(4-tetrahydropyranyl)ammonium chloride, N,N-dimethyl-N-(((7-(4-methylphenyl)-2,3-dihydro-1-benzoxepin-4-yl)carbonyl)amino)benzyl)-N-(4-oxocyclohexyl)ammonium chloride, N-(4-(((7-(4-ethoxyphenyl)-2,3-dihydro-1-benzoxepin-4-yl)carbonyl)amino)benzyl)-N,N-dimethyl-N-(4-tetrahydropyranyl)ammonium chloride, N-[4-[N-methyl-N-(tetrahydropyran-4-yl)aminomethyl]phenyl]-7-(4-propoxyphenyl)-1,1-dioxo-2,3-dihydro-1-benzothiepin-4-carboxamide, 7-(4-butoxyphenyl)-N-[4-[N-methyl-N-(tetrahydropyran-4-yl)aminomethyl]phenyl]-1,1-dioxo-2,3-dihydro-1-benzothiepin-4-carboxamide, 7-[4-[N-methyl-N-(2-propoxyethyl)amino]phenyl]-N-[4-[N-methyl-N-(tetrahydropyran-4-yl)amino]methyl]phenyl]-1,1-dioxo-2,3-dihydro-1-benzothiepin-4-carboxamide, 7-[4-(2-ethoxyethoxy)phenyl]-N-[4-[N-methyl-N-(tetrahydropyran-4-yl)amino]methyl]phenyl]-1,1-dioxo-2,3-dihydro-1-benzothiepin-4-carboxamide, N-[4-[N-methyl-N-(tetrahydropyran-4-yl)amino]methyl]phenyl]-7-[4-(2-propoxyethoxy)phenyl]-1,1-dioxo-2,3-dihydro-1-benzothiepin-4-carboxamide, 7-[4-(2-butoxyethoxy)phenyl]-N-[4-[N-methyl-N-(tetrahydropyran-4-yl)amino]methyl]phenyl]-1,1-dioxo-2,3-dihydro-1-benzothiepin-4-carboxamide, 7-[4-(2-ethoxyethoxy)-3,5-dimethylphenyl]-N-[4-[N-methyl-N-(tetrahydro-2H-pyran-4-yl)amino]methyl]phenyl]-1,1-dioxo-2,3-dihydro-1-benzothiepin-4-carboxamide, 7-[2-chloro-4-(2-propoxyethoxy)phenyl]-N-[4-[N-methyl-N-(tetrahydropyran-4-yl)amino]methyl]phenyl]-1,1-dioxo-2,3-dihydro-1-benzothiepin-4-carboxamide, 7-(3-methyl-4-propoxyphenyl)-N-[4-[N-methyl-N-(tetrahydropyran-4-yl)amino]methyl]phenyl]-1,1-dioxo-2,3-dihydro-1-benzothiepin-4-carboxamide, 7-(3,4-dipropoxyphenyl)-N-(4-((N-methyl-N-(tetrahydro-2H-pyran-4-yl)amino)methyl)phenyl)-1,1-dioxo-2,3-dihydro-1-benzothiepin-4-carboxamide, 7-[4-(2-

ethoxyethoxy)phenyl]-1-ethyl-N-[4-[[N-methyl-N-(tetrahydropyran-4-yl)amino]methyl]phenyl]-2,3-dihydro-1-benzazepin-4-carboxamide, 1-ethyl-7-[4-(2-propoxyethoxy)phenyl]-N-[4-[[N-methyl-N-(tetrahydropyran-4-yl)amino]methyl]phenyl]-2,3-dihydro-1-benzazepin-4-carboxamide, 7-[4-(2-butoxyethoxy)phenyl]-1-ethyl-N-[4-[[N-methyl-N-(tetrahydropyran-4-yl)amino]methyl]phenyl]-2,3-dihydro-1-benzazepin-4-carboxamide, 7-[4-(2-ethoxyethoxy)phenyl]-1-formyl-N-[4-[[N-methyl-N-(tetrahydropyran-4-yl)amino]methyl]phenyl]-2,3-dihydro-1-benzazepin-4-carboxamide, 1-formyl-7-[4-(2-propoxyethoxy)phenyl]-N-[4-[[N-methyl-N-(tetrahydropyran-4-yl)amino]methyl]phenyl]-2,3-dihydro-1-benzazepin-4-carboxamide, 7-[4-(2-butoxyethoxy)phenyl]-1-formyl-N-[4-[[N-methyl-N-(tetrahydropyran-4-yl)amino]methyl]phenyl]-2,3-dihydro-1-benzazepin-4-carboxamide, 7-[4-(2-butoxyethoxy)phenyl]-N-[4-[[N-methyl-N-(tetrahydropyran-5-yl)amino]methyl]phenyl]-1-propyl-2,3-dihydro-1-benzazepin-4-carboxamide, N-[4-[[N-methyl-N-(tetrahydropyran-5-yl)amino]methyl]phenyl]-7-[4-(2-propoxyethoxy)phenyl]-1-propyl-2,3-dihydro-1-benzazepin-4-carboxamide, 1-benzyl-7-[4-(2-butoxyethoxy)phenyl]-N-[4-[[N-methyl-N-(tetrahydropyran-4-yl)amino]methyl]phenyl]-2,3-dihydro-1-benzazepin-4-carboxamide, 7-[4-(2-butoxyethoxy)phenyl]-1-cyclopropylmethyl-N-[4-[[N-methyl-N-(tetrahydropyran-4-yl)amino]methyl]phenyl]-2,3-dihydro-1-benzazepin-4-carboxamide, 7-[4-(2-butoxyethoxy)phenyl]-N-[4-[[N-methyl-N-(tetrahydropyran-4-yl)amino]methyl]phenyl]-1-phenyl-2,3-dihydro-1-benzazepin-4-carboxamide, 7-[4-(2-butoxyethoxy)phenyl]-1-(3,4-methylenedioxy)phenyl-N-[4-[[N-methyl-N-(tetrahydropyran-4-yl)amino]methyl]phenyl]-2,3-dihydro-1-benzazepin-4-carboxamide, 7-[4-(2-butoxyethoxy)phenyl]-1-(2-methyloxazol-5-yl)-N-[4-[[N-methyl-N-(tetrahydropyran-4-yl)amino]methyl]phenyl]-2,3-dihydro-1-benzazepin-4-carboxamide, 1-allyl-7-[4-(2-butoxyethoxy)phenyl]-N-[4-[[N-methyl-N-(tetrahydropyran-4-yl)amino]methyl]phenyl]-2,3-dihydro-1-benzazepin-4-carboxamide, 7-[4-(2-butoxyethoxy)phenyl]-N-[4-[[N-methyl-N-(tetrahydropyran-4-yl)amino]methyl]phenyl]-1-(3-thienyl)methyl-2,3-dihydro-1-benzazepin-4-carboxamide, 7-[4-(2-butoxyethoxy)phenyl]-N-[4-[[N-methyl-N-(tetrahydropyran-4-yl)amino]methyl]phenyl]-1-(thiazol-2-yl)methyl-2,3-dihydro-1-benzazepin-4-carboxamide, 7-[4-(2-butoxyethoxy)phenyl]-1-(1-methylpyrazol-4-yl)methyl-N-[4-[[N-methyl-N-(tetrahydropyran-4-yl)amino]methyl]phenyl]-2,3-dihydro-1-benzazepin-4-carboxamide, 7-[4-(2-butoxyethoxy)phenyl]-1-(3-methylisothiazol-5-yl)methyl-N-[4-[[N-

methyl-N-(tetrahydropyran-4-yl)amino]methyl]phenyl]-2,3-dihydro-1-benzoazepin-4-carboxamide, 7-[4-(2-butoxyethoxy)phenyl]-1-(1-ethylpyrazol-4-yl)methyl-N-[4-[[N-methyl-N-(tetrahydropyran-4-yl)amino]methyl]phenyl]-2,3-dihydro-1-benzoazepin-4-carboxamide, 7-[4-(2-butoxyethoxy)phenyl]-1-isobutyl-N-[4-[[N-methyl-N-(tetrahydropyran-5-yl)amino]methyl]phenyl]-2,3-dihydro-1-benzoazepin-4-carboxamide, 1-isobutyl-N-[4-[[N-methyl-N-(tetrahydropyran-5-yl)amino]methyl]phenyl]-7-[4-(2-propoxyethoxy)phenyl]-2,3-dihydro-1-benzoazepin-4-carboxamide, 7-[4-(2-butoxyethoxy)phenyl]-N-[4-[[N-methyl-N-(tetrahydropyran-4-yl)amino]methyl]phenyl]-1-(thiazol-5-yl)methyl-2,3-dihydro-1-benzoazepin-4-carboxamide, 7-[4-(2-butoxyethoxy)phenyl]-N-[4-[[N-methyl-N-(tetrahydropyran-4-yl)amino]methyl]phenyl]-1-(1-methyltetrazol-5-yl)methyl-2,3-dihydro-1-benzoazepin-4-carboxamide, or 7-[4-(2-butoxyethoxy)phenyl]-N-[4-[[N-methyl-N-(tetrahydropyran-4-yl)amino]methyl]phenyl]-1-(2-methyltetrazol-5-yl)methyl-2,3-dihydro-1-benzoazepin-4-carboxamide or a salt thereof.

5. (Cancelled)

6. (Withdrawn) A method of preventing or treating chronic rheumatoid arthritis, autoimmune diseases, allergic disorders, ischemic brain cell damage, myocardial infarction, chronic nephritis, and arteriosclerosis, which comprises the step of administering an effective amount of a compound having a CCR antagonist effect to a mammal.

7. (Cancelled)

8. (Withdrawn) Use of a compound having a CCR antagonist effect for manufacturing an agent for the prevention or treatment of chronic rheumatoid arthritis, autoimmune diseases, allergic disorders, ischemic brain cell damage, myocardial infarction, chronic nephritis, and arteriosclerosis.

9-11. (Cancelled)

12. (Currently amended) The method of ~~claim 5~~claim 1, wherein the compound having a CCR antagonist effect is 1-Acetyl-N-(3-{4-[4-(aminocarbonyl)benzyl]-1-piperidinyl}propyl)-N-(3-chloro-4-methylphenyl)-4-piperidinecarboxamide, or a salt thereof.